

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121522\
 Data File : BG055978.D
 Acq On : 15 Dec 2022 15:44
 Operator : CG/JU
 Sample : N6057-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1521

Quant Time: Dec 16 04:03:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 14 02:24:06 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.376	152	6831	20.000	ng	0.00
21) Naphthalene-d8	11.208	136	26099	20.000	ng	# 0.00
39) Acenaphthene-d10	14.980	164	23658	20.000	ng	0.00
64) Phenanthrene-d10	17.718	188	67714	20.000	ng	# 0.00
76) Chrysene-d12	22.024	240	71930	20.000	ng	0.00
86) Perylene-d12	25.514	264	87457	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.890	112	31408	107.828	ng	0.00
7) Phenol-d6	7.500	99	40148	107.169	ng	0.00
23) Nitrobenzene-d5	9.545	82	25505	68.292	ng	0.00
42) 2,4,6-Tribromophenol	16.454	330	60723	106.445	ng	0.00
45) 2-Fluorobiphenyl	13.611	172	102659	60.858	ng	0.00
79) Terphenyl-d14	20.291	244	266483	65.478	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.753	178	9606	2.707	ng	# 94
75) Fluoranthene	19.745	202	19978	4.020	ng	98
78) Pyrene	20.103	202	14595	3.353	ng	97
83) Chrysene	22.071	228	9595	2.023	ng	# 91
88) Benzo(b)fluoranthene	24.392	252	12842	2.269	ng	# 98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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